

=> file registry

FILE 'REGISTRY' ENTERED AT 15:12:34 ON 18 JUL 2005

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STRUCTURE FILE UPDATES: 17 JUL 2005 HIGHEST RN 855596-49-5

DICTIONARY FILE UPDATES: 17 JUL 2005 HIGHEST RN 855596-49-5

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH JANUARY 18, 2005

Please note that search-term pricing does apply when conducting SmartSELECT searches.

\*\*\*\*\*  
\*  
\* The CA roles and document type information have been removed from \*  
\* the IDE default display format and the ED field has been added, \*  
\* effective March 20, 2005. A new display format, IDERL, is now \*  
\* available and contains the CA role and document type information. \*  
\*  
\*\*\*\*\*

Structure search iteration limits have been increased. See HELP SLIMITS for details.

Experimental and calculated property data are now available. For more information enter HELP PROP at an arrow prompt in the file or refer to the file summary sheet on the web at:  
<http://www.cas.org/ONLINE/DBSS/registryss.html>

=> file caplus

FILE 'CAPLUS' ENTERED AT 15:12:42 ON 18 JUL 2005

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FILE COVERS 1907 - 18 Jul 2005 VOL 143 ISS 4

FILE LAST UPDATED: 17 Jul 2005 (20050717/ED)

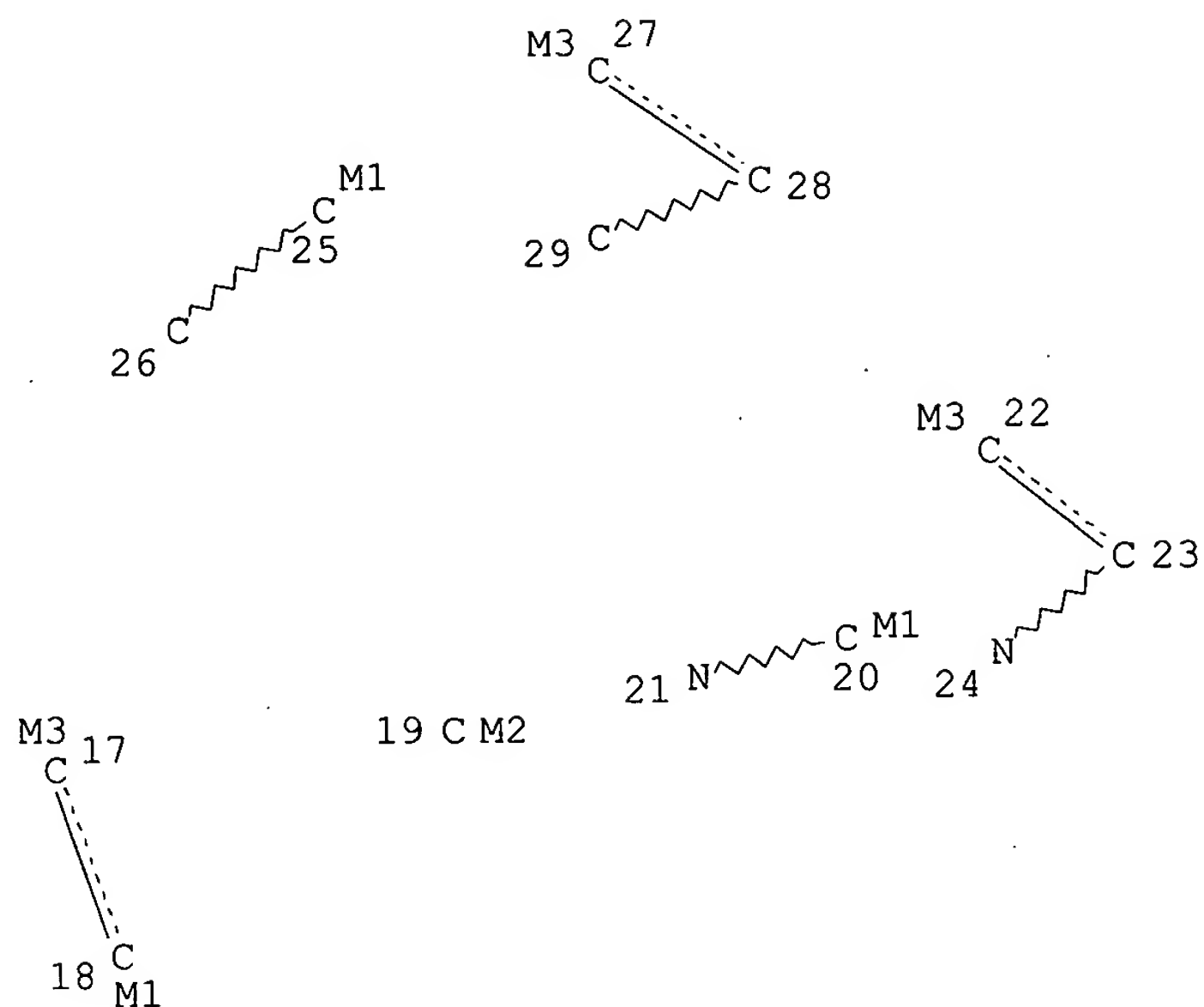
New CAS Information Use Policies, enter HELP USAGETERMS for details.

This file contains CAS Registry Numbers for easy and accurate

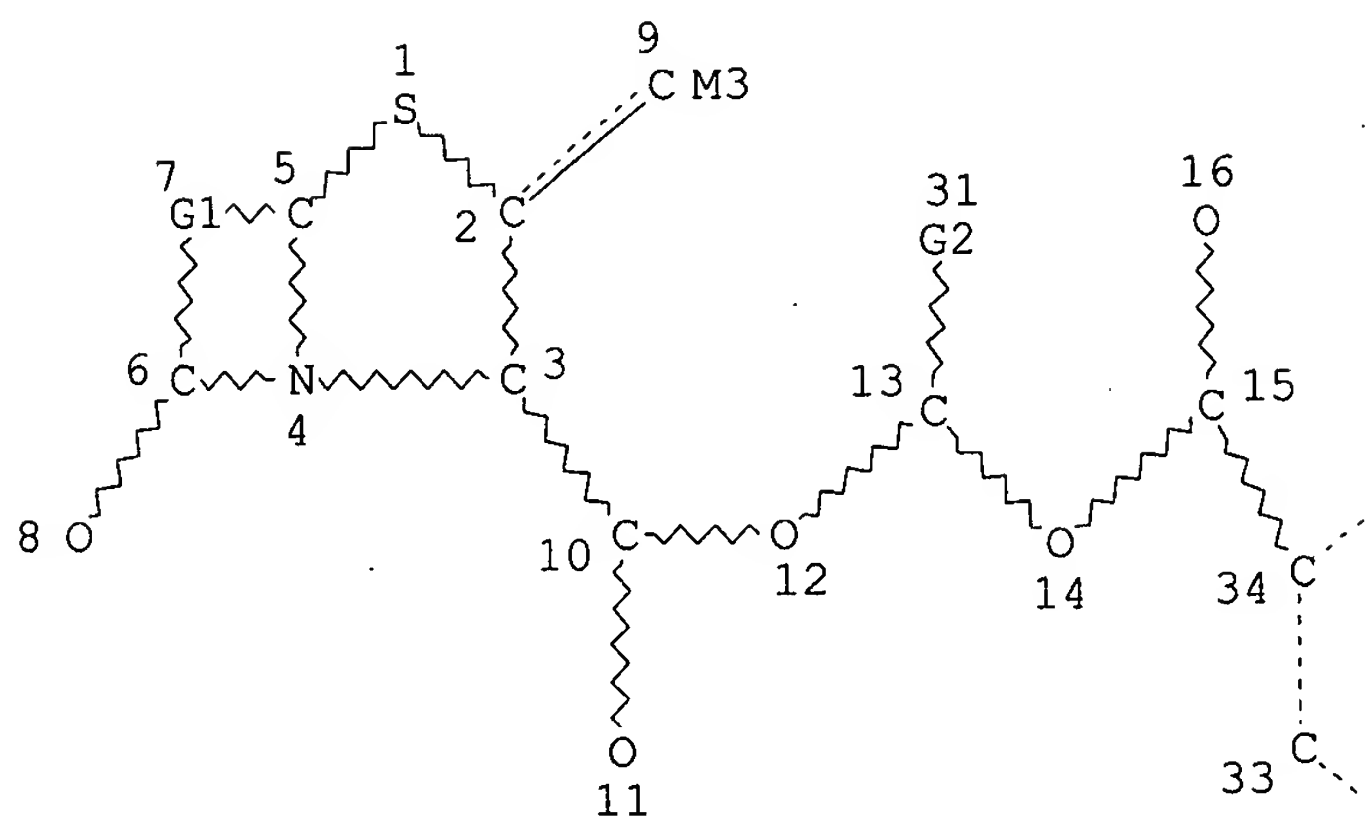
substance identification.

=> d stat que L13  
L3 STR

H 38

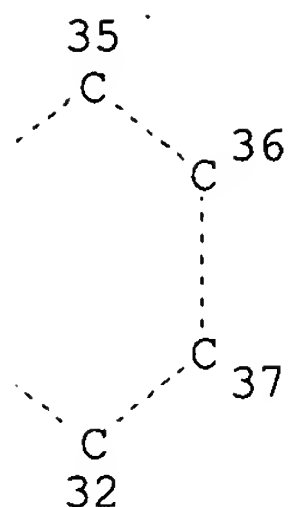


Page 1-A



Ak 30

Page 2-A



Page 2-B

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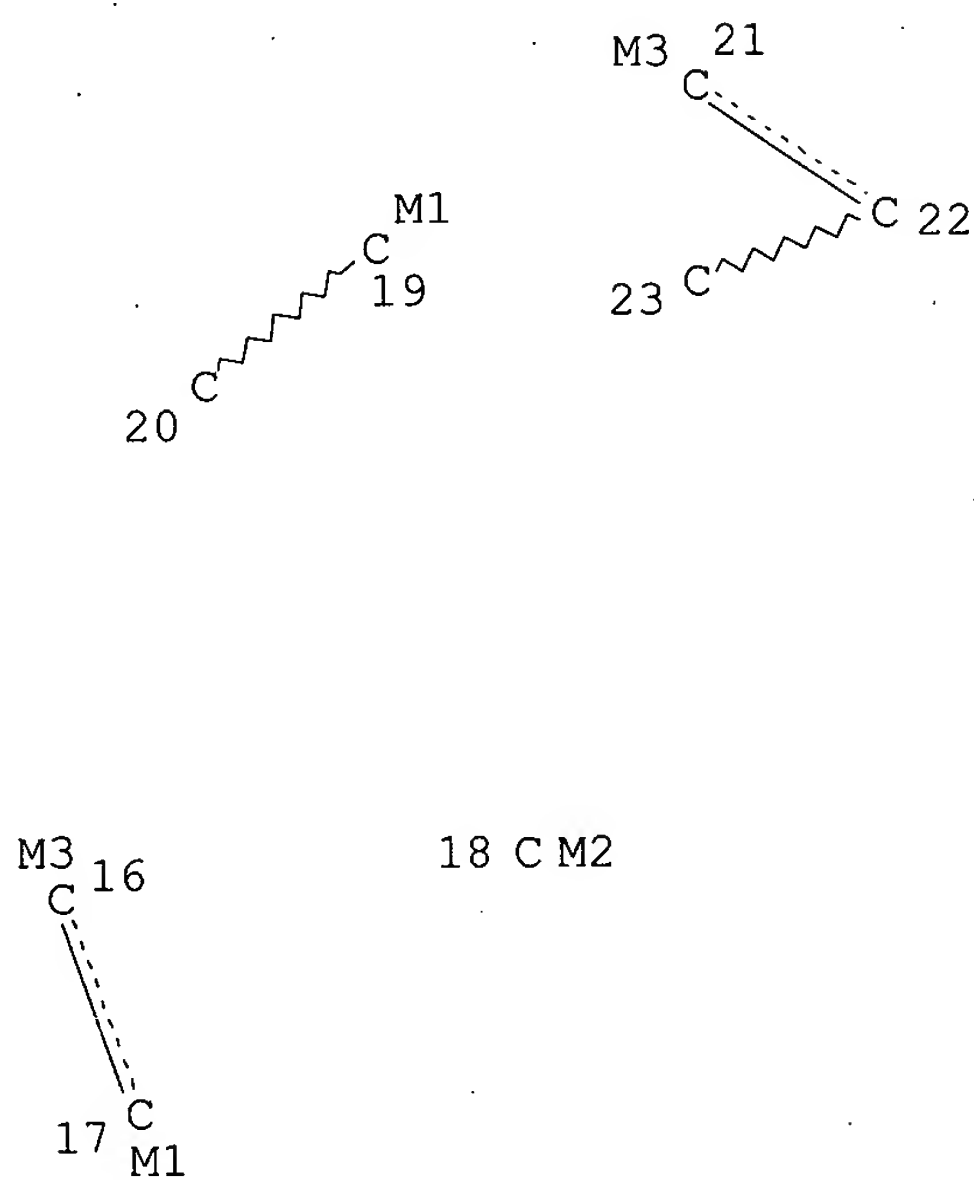
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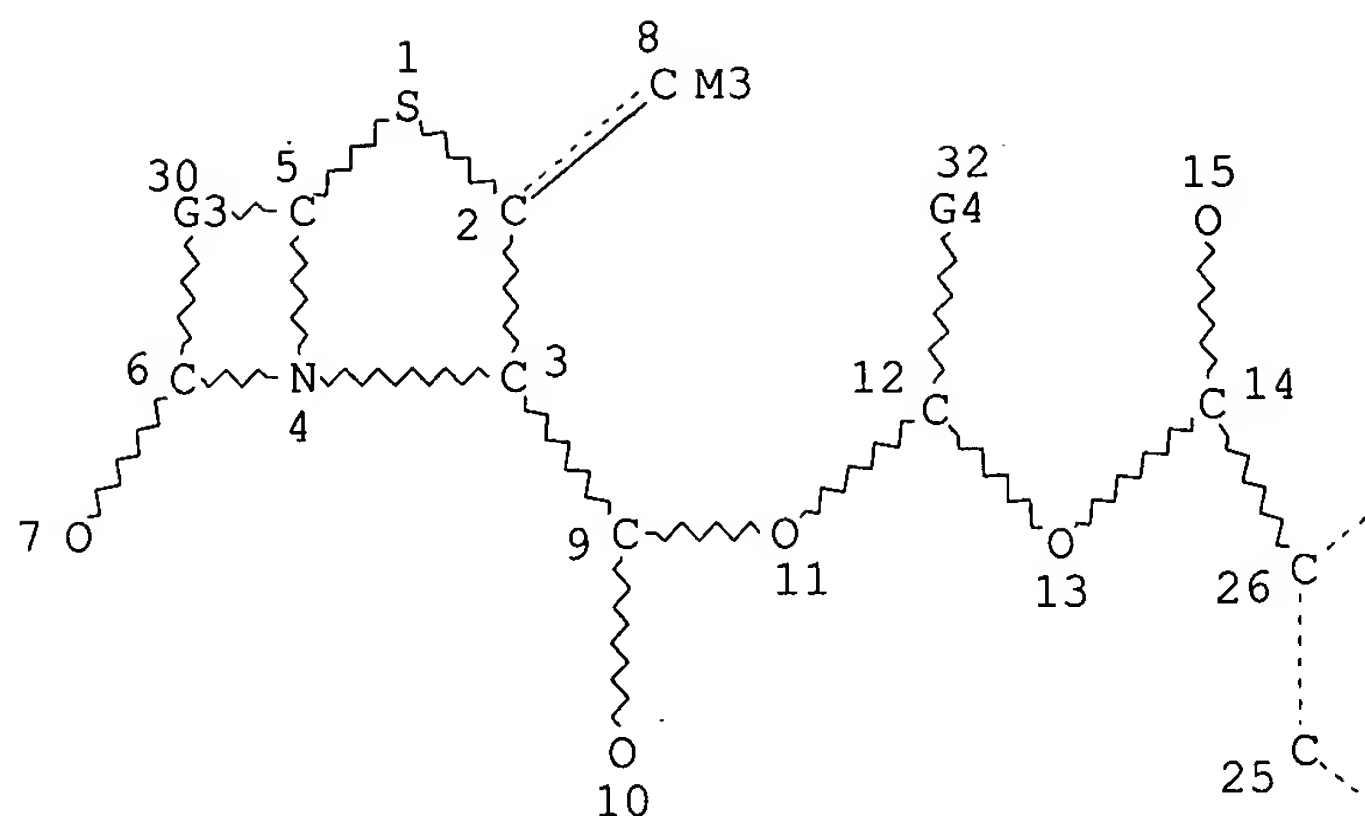
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GRAPH ATTRIBUTES:  
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NUMBER OF NODES IS 38

STEREO ATTRIBUTES: NONE  
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L8 STR

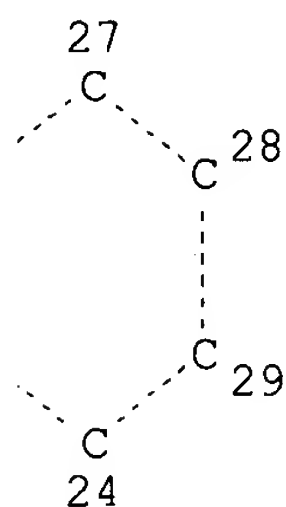
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Page 2-B

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VAR G4=33/31

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DEFAULT ECLEVEL IS LIMITED  
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GRAPH ATTRIBUTES:  
RING(S) ARE ISOLATED OR EMBEDDED  
NUMBER OF NODES IS 33

STEREO ATTRIBUTES: NONE  
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L13 4 SEA FILE=CAPLUS ABB=ON PLU=ON L12

=> d ibib abs hitstr L13 1-4

L13 ANSWER 1 OF 4 CAPLUS COPYRIGHT 2005 ACS on STN  
ACCESSION NUMBER: 2004:182891 CAPLUS  
DOCUMENT NUMBER: 140:217438  
TITLE: Preparation of hydroxymethylpenicillanic acid sulfones  
as  $\beta$ -lactamase inhibitor prodrugs  
INVENTOR(S): Marfat, Anthony; McLeod, Dale Gordon  
PATENT ASSIGNEE(S): Pfizer Products Inc., USA  
SOURCE: PCT Int. Appl., 83 pp.  
CODEN: PIXXD2  
DOCUMENT TYPE: Patent  
LANGUAGE: English  
FAMILY ACC. NUM. COUNT: 1  
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2004018484	A1	20040304	WO 2003-IB3582	20030811

W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NI, NO, NZ, OM, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW

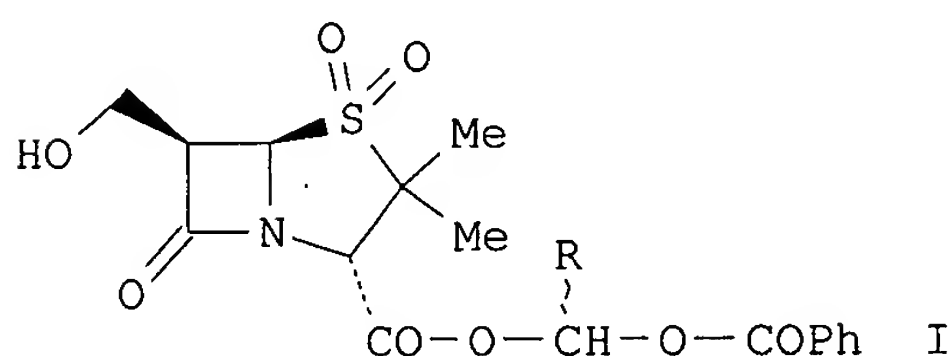
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CA 2494953 AA 20040304 CA 2003-2494953 20030811  
 EP 1534717 A1 20050601 EP 2003-792574 20030811

R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, MK, CY, AL, TR, BG, CZ, EE, HU, SK

US 2004110740 A1 20040610 US 2003-648408 20030825  
 PRIORITY APPLN. INFO.: US 2002-405640P P 20020823  
 WO 2003-IB3582 W 20030811

GI



AB Prodrugs of 6β-hydroxymethylpenicillanic acid sulfone of formula I [R = H, Me] and solvates thereof, are prepared. Also disclosed are pharmaceutical compns. comprising a prodrug of the present invention, or a solvate thereof, an optional β-lactam antibiotic and at least one pharmaceutically acceptable carrier. Further disclosed is a method for increasing the therapeutic effectiveness of a β-lactam antibiotic in a mammal by administering an effective amount of a β-lactam antibiotic and an effectiveness-increasing amount of a prodrug of the present invention, or a solvate thereof. Addnl. disclosed is a method for treating a bacterial infection in a mammal by administering a therapeutically effective amount of a pharmaceutical composition of the present invention to a mammal in need thereof. Thus, a combination of amoxicillin and prodrug I (R = (R)-Me) was effective against *S. pneumoniae* in gerbil otitis media model.

IT 666174-85-2P 666174-86-3P 666174-87-4P  
 666174-88-5P

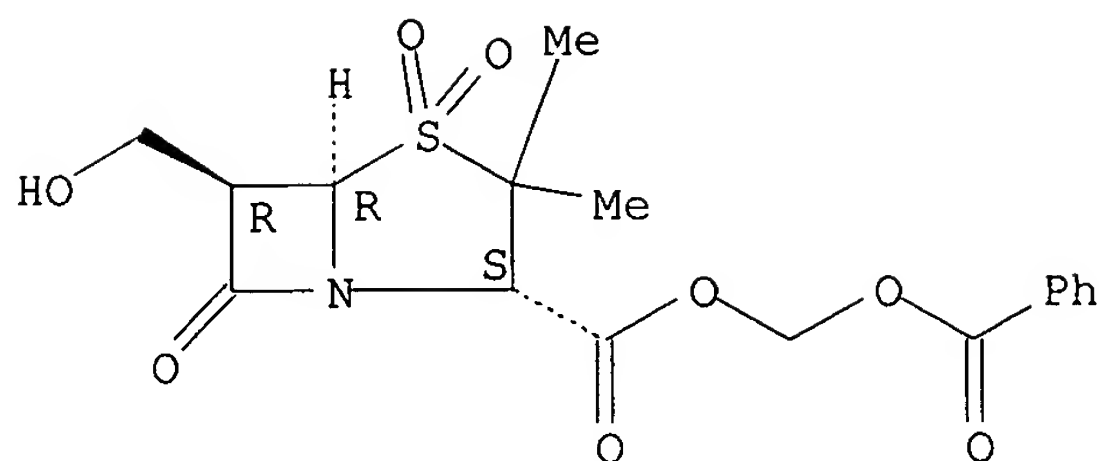
RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(preparation of hydroxymethylpenicillanic acid sulfones as β-lactamase inhibitor prodrugs)

RN 666174-85-2 CAPLUS

CN 4-Thia-1-azabicyclo[3.2.0]heptane-2-carboxylic acid, 6-(hydroxymethyl)-3,3-dimethyl-7-oxo-, (benzoyloxy)methyl ester, 4,4-dioxide, (2S,5R,6R)- (9CI)  
 (CA INDEX NAME)

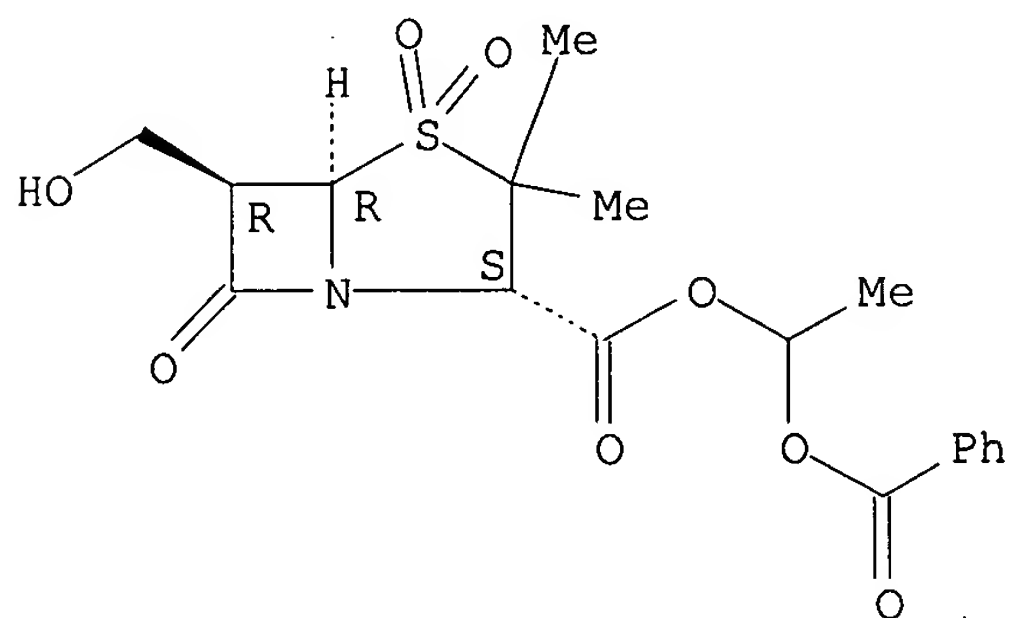
Absolute stereochemistry.



RN 666174-86-3 CAPLUS

CN 4-Thia-1-azabicyclo[3.2.0]heptane-2-carboxylic acid, 6-(hydroxymethyl)-3,3-dimethyl-7-oxo-, 1-(benzoyloxy)ethyl ester, 4,4-dioxide, (2S,5R,6R)- (9CI)  
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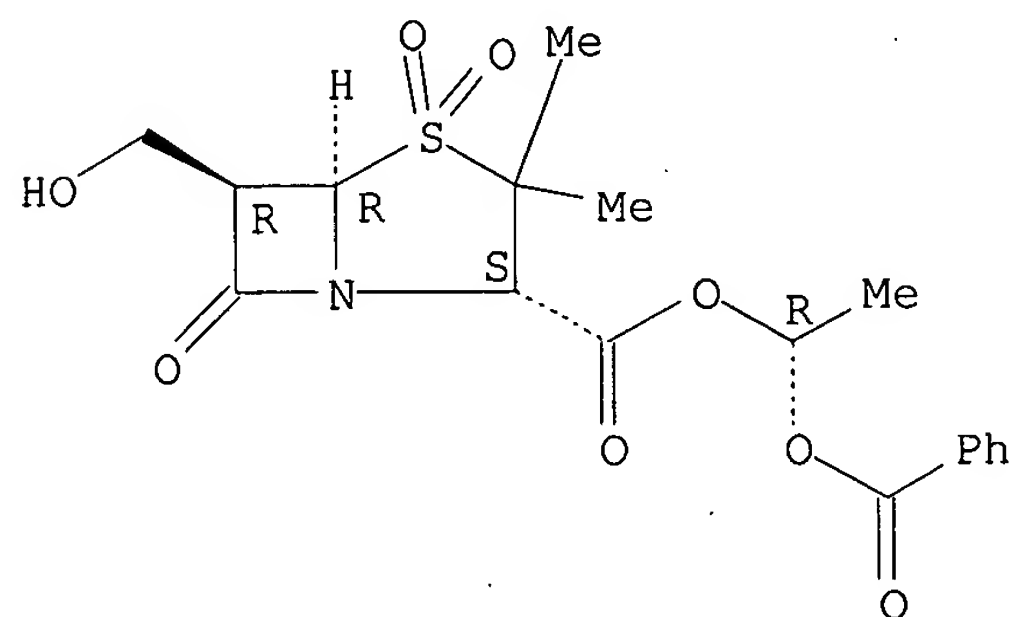
Absolute stereochemistry.



RN 666174-87-4 CAPLUS

CN 4-Thia-1-azabicyclo[3.2.0]heptane-2-carboxylic acid, 6-(hydroxymethyl)-3,3-dimethyl-7-oxo-, (1R)-1-(benzoyloxy)ethyl ester, 4,4-dioxide, (2S,5R,6R)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (+).

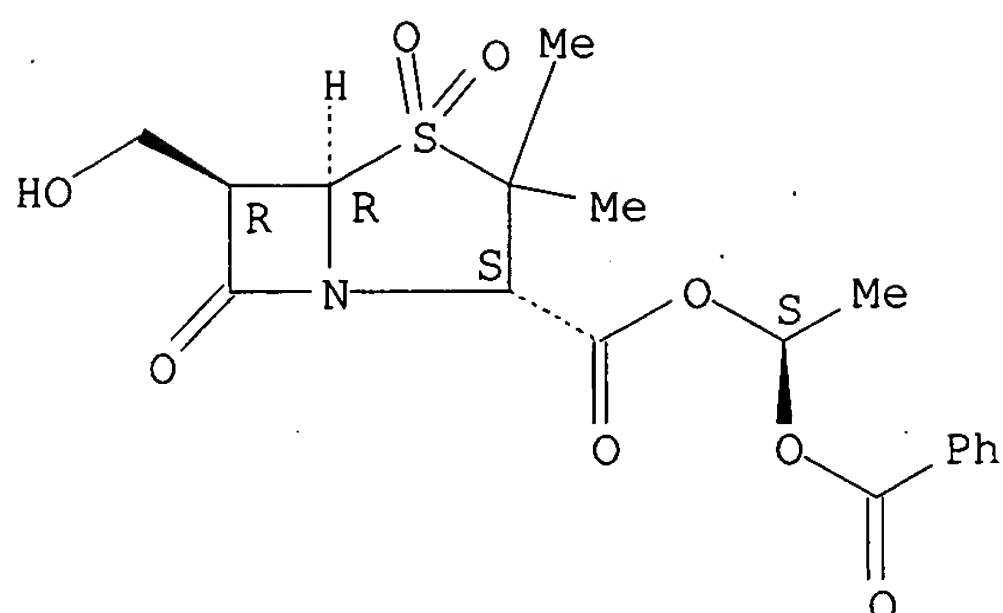


RN 666174-88-5 CAPLUS

CN 4-Thia-1-azabicyclo[3.2.0]heptane-2-carboxylic acid, 6-(hydroxymethyl)-3,3-dimethyl-7-oxo-, (1S)-1-(benzoyloxy)ethyl ester, 4,4-dioxide, (2S,5R,6R)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.





REFERENCE COUNT: 5 THERE ARE 5 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L13 ANSWER 2 OF 4 CAPLUS COPYRIGHT 2005 ACS on STN

ACCESSION NUMBER: 1984:591536 CAPLUS

DOCUMENT NUMBER: 101:191536

TITLE: 1,1-Alkanediol dicarboxylate-linked antibacterial agents

INVENTOR(S): Jasys, Vytautas J.; Kellogg, Michael S.

PATENT ASSIGNEE(S): Pfizer Inc., USA

SOURCE: U.S., 39 pp. Cont.-in-part of U.S. Ser. No. 334,022, abandoned.

CODEN: USXXAM

DOCUMENT TYPE: Patent

LANGUAGE: English

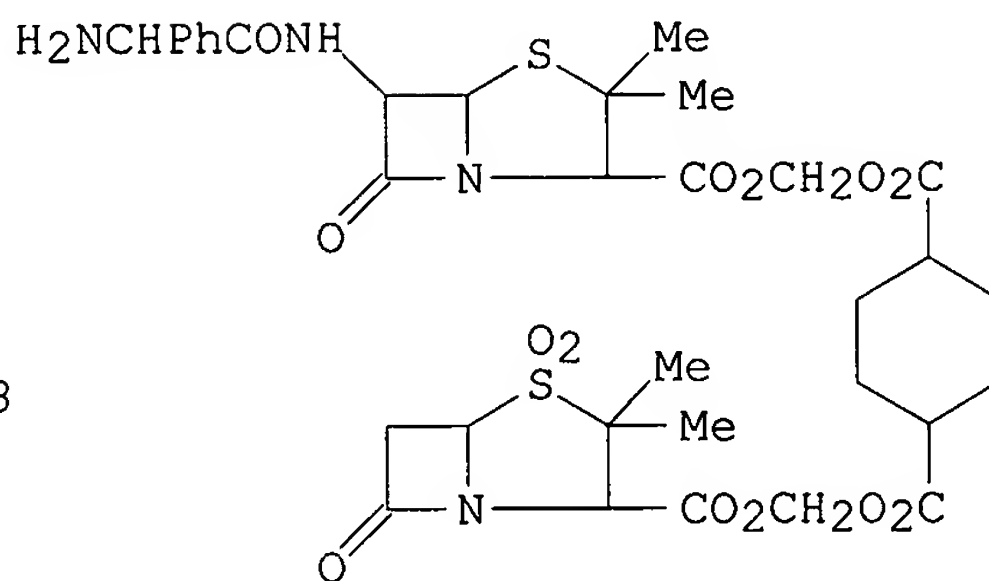
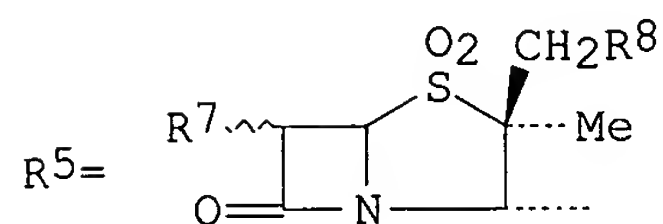
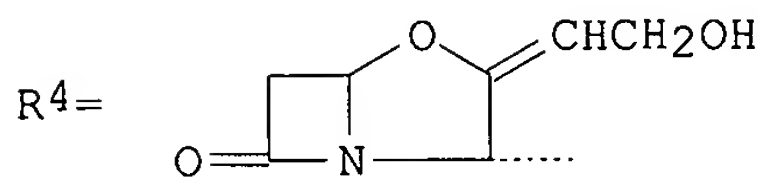
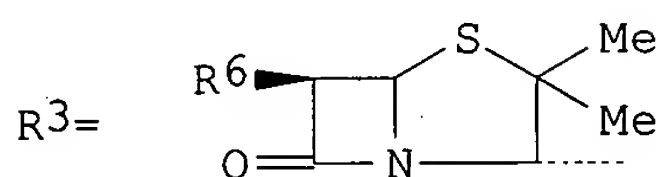
FAMILY ACC. NUM. COUNT: 2

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
US 4457924	A	19840703	US 1982-429915	19820930
EP 83484	A1	19830713	EP 1982-306683	19821214
EP 83484	B1	19860219		
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RO 87709	B3	19851031	RO 1982-113244	19821220
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FI 8204409	A	19830623	FI 1982-4409	19821221
FI 80039	B	19891229		
FI 80039	C	19900410		
NO 8204305	A	19830623	NO 1982-4305	19821221
AU 8291721	A1	19830630	AU 1982-91721	19821221
AU 537214	B2	19840614		
ZA 8209372	A	19830928	ZA 1982-9372	19821221
HU 27683	O	19831028	HU 1982-4105	19821221
HU 187737	B	19860228		
ES 518425	A1	19840201	ES 1982-518425	19821221
DD 207379	A5	19840229	DD 1982-246325	19821221
IL 67530	A1	19860228	IL 1982-67530	19821221
CA 1213582	A1	19861104	CA 1982-418192	19821221
PL 140291	B1	19870430	PL 1982-248637	19821221
PL 141306	B1	19870731	PL 1982-239651	19821221
SU 1405704	A3	19880623	SU 1982-3529507	19821221

PL 145927	B1	19881130	PL 1982-256903	19821221
JP 58116486	A2	19830711	JP 1982-225773	19821222
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CS 236867	B2	19850515	CS 1982-9559	19821222
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FI 81353	B	19900629		
FI 81353	C	19901010		
JP 02270881	A2	19901105	JP 1990-33601	19900214
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GI



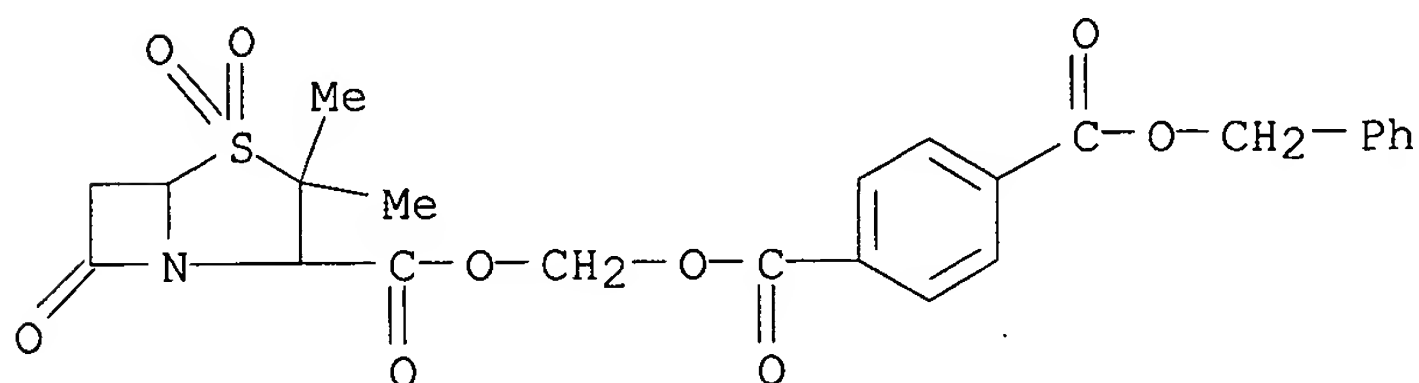
AB RCO2CHR1O2CXO2C(CHR1O2C)nR2 [X = C1-12 alkylene, alkylidene (un)substituted by Ph or CO2H, cycloalkylene, phenylene, naphthalenediyl, furandiyl, thiophendiyl, pyridinediyl, pyrazinediyl; R = R3-R5; R1 = H, alkyl; R2 = R3-R5, H, alkyl, CH2Ph, CHR1Cl, CHR1I, NBu4; R6 = NH2, 2,6-(MeO)2C6H3CONH, PhOCH2CONH, 4-R9C6H4CHR1OCONH; R7 = H, CH2OH, CH2NH2, CHMeNH2; R8 = H, Cl, OAc; R9 = H, OH, acyloxy, alkoxy carbonyloxy, (un)substituted BzO; R10 = H, (un)protected NH2, N3] were prepared Thus, I was prepared from Na penicillanate 1,1-dioxide, ampicillin, K benzyl trans-1,4-cyclohexanedicarboxylate, ClCH2I, and ClCH2Br in 10 steps.

IT **87343-43-9P 87343-50-8P**  
 RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(preparation and hydrogenolysis of)

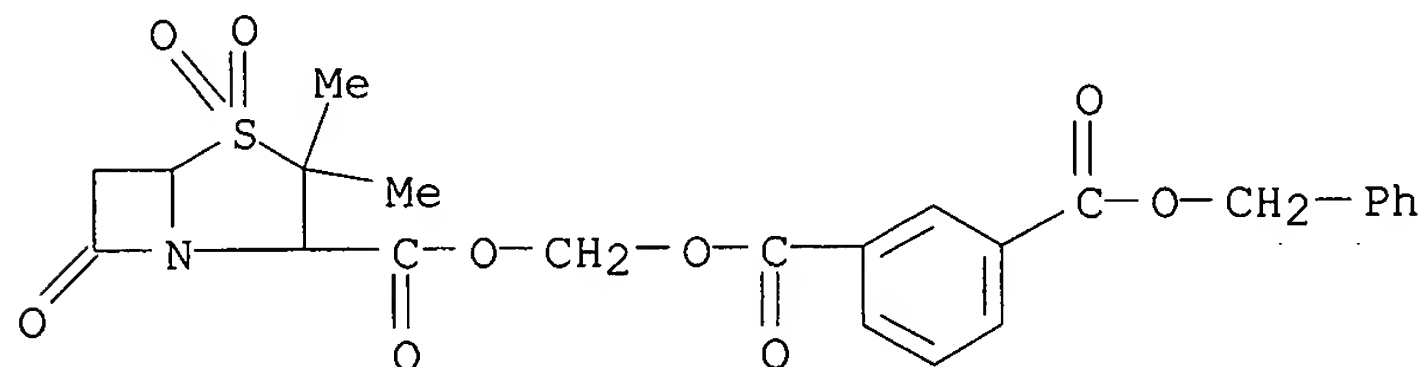
RN 87343-43-9 CAPLUS

CN 1,4-Benzenedicarboxylic acid, [[[3,3-dimethyl-4,4-dioxido-7-oxo-4-thia-1-azabicyclo[3.2.0]hept-2-yl)carbonyl]oxy]methyl phenylmethyl ester, (2S-cis)- (9CI) (CA INDEX NAME)



RN 87343-50-8 CAPLUS

CN 1,3-Benzenedicarboxylic acid, [[[3,3-dimethyl-4,4-dioxido-7-oxo-4-thia-1-azabicyclo[3.2.0]hept-2-yl)carbonyl]oxy]methyl phenylmethyl ester, (2S-cis)- (9CI) (CA INDEX NAME)



IT 87352-91-8P 87352-93-0P

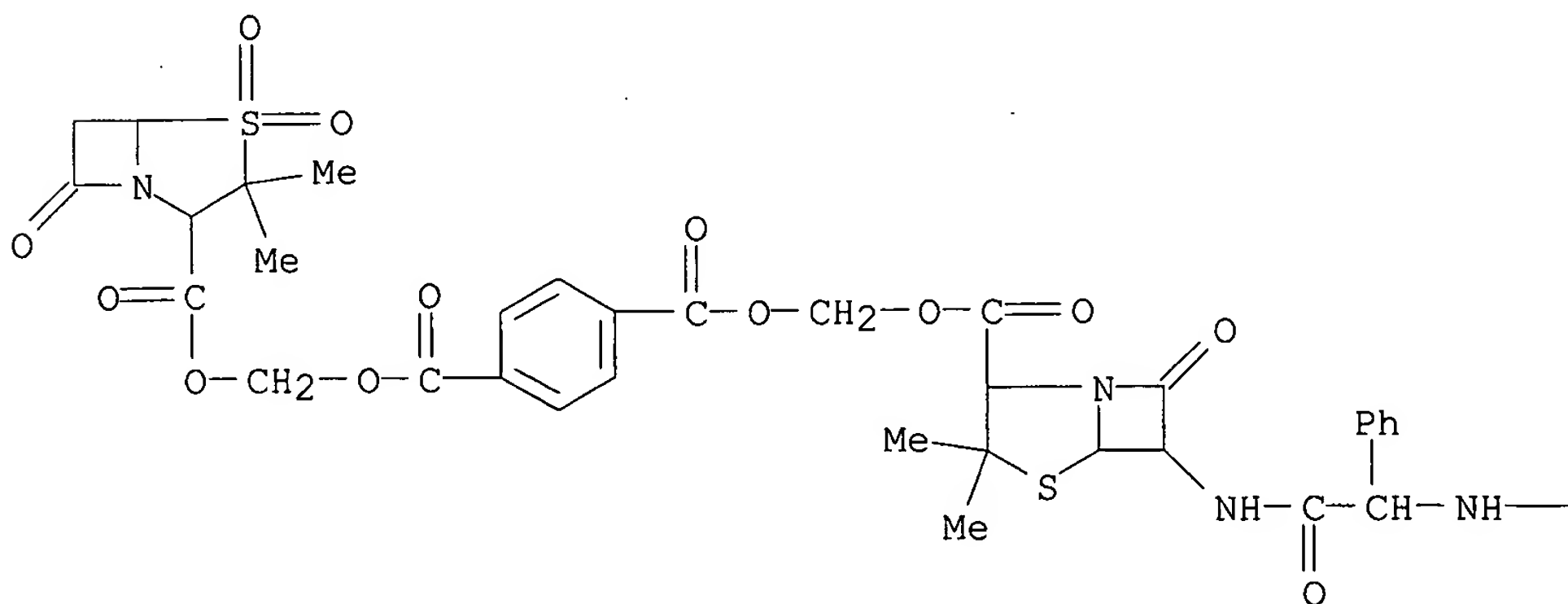
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(preparation and hydrolysis of)

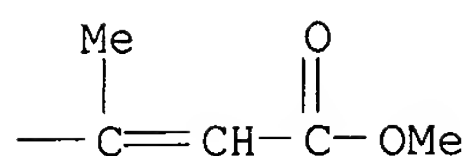
RN 87352-91-8 CAPLUS

CN 1,4-Benzenedicarboxylic acid, [[[3,3-dimethyl-4,4-dioxido-7-oxo-4-thia-1-azabicyclo[3.2.0]hept-2-yl)carbonyl]oxy]methyl [[[6-[[[(3-methoxy-1-methyl-3-oxo-1-propenyl)amino]phenylacetyl]amino]-3,3-dimethyl-7-oxo-4-thia-1-azabicyclo[3.2.0]hept-2-yl)carbonyl]oxy]methyl ester, [2S-[2α(2R\*,5S\*),5α,6β(S\*)]]- (9CI) (CA INDEX NAME)

PAGE 1-A



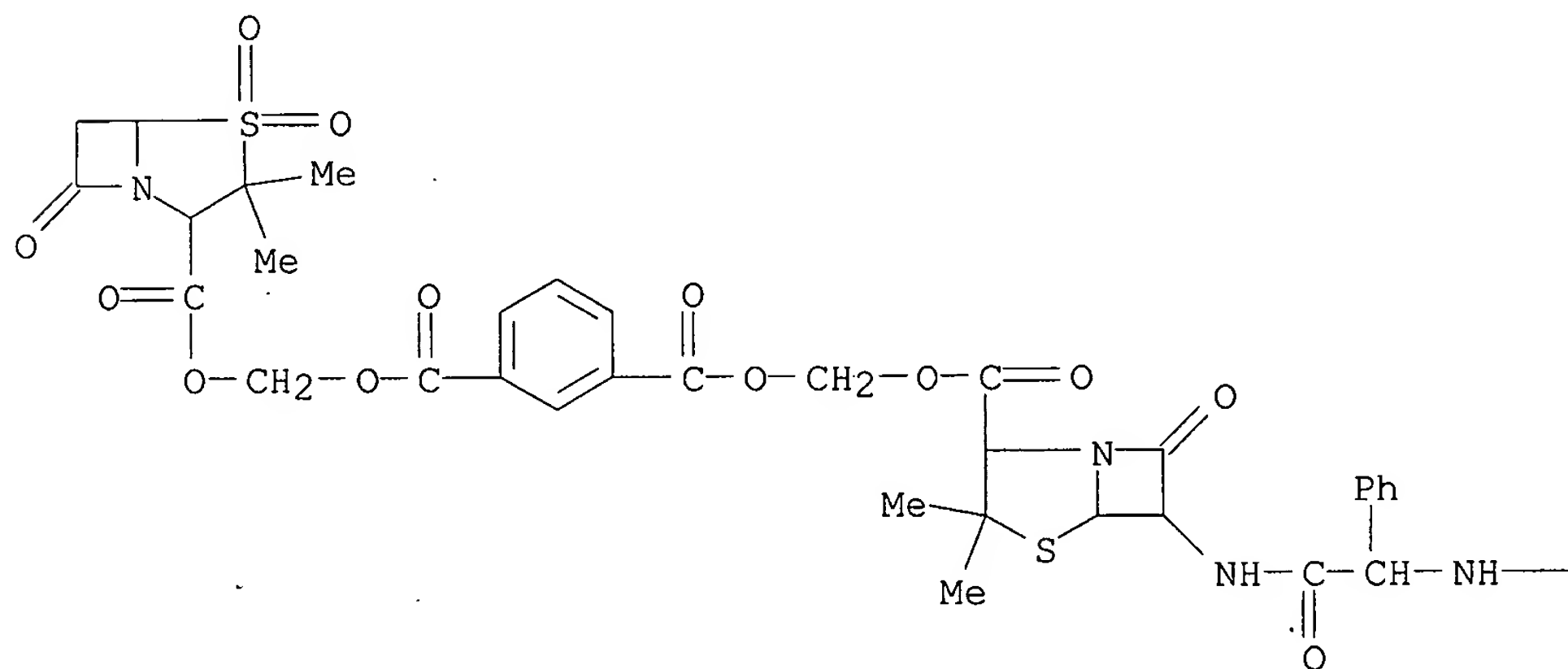
PAGE 1-B



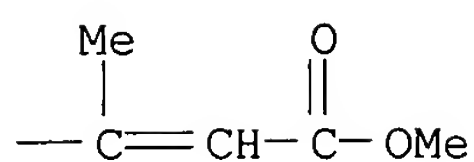
RN 87352-93-0 CAPLUS

CN 1,3-Benzenedicarboxylic acid, [[(3,3-dimethyl-4,4-dioxido-7-oxo-4-thia-1-azabicyclo[3.2.0]hept-2-yl)carbonyl]oxy]methyl [[[6-[[[(3-methoxy-1-methyl-3-oxo-1-propenyl)amino]phenylacetyl]amino]-3,3-dimethyl-7-oxo-4-thia-1-azabicyclo[3.2.0]hept-2-yl]carbonyl]oxy]methyl ester, [2S-[2 $\alpha$ (2R\*,5S\*),5 $\alpha$ ,6 $\beta$ (S\*)]]- (9CI) (CA INDEX NAME)

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PAGE 1-B



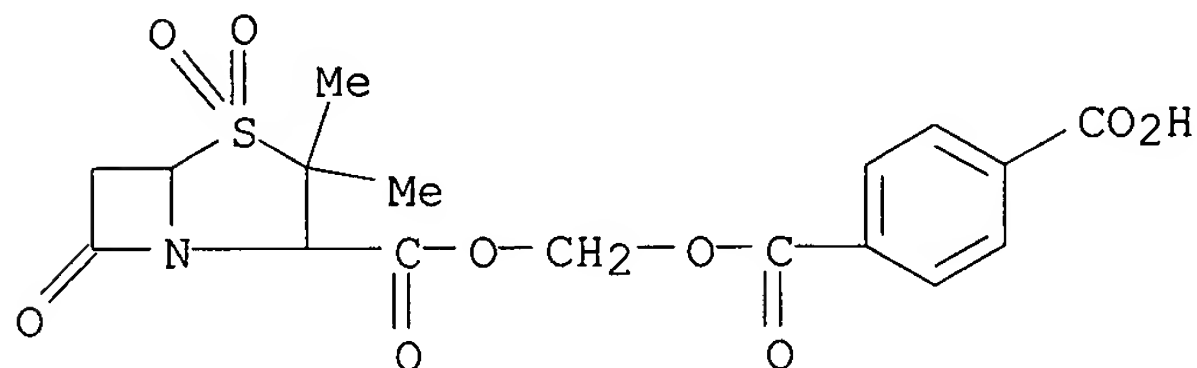
IT 87343-44-0P

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(preparation and neutralization of)

RN 87343-44-0 CAPLUS

CN 1,4-Benzenedicarboxylic acid, mono[[[(3,3-dimethyl-4,4-dioxido-7-oxo-4-thia-1-azabicyclo[3.2.0]hept-2-yl)carbonyl]oxy)methyl] ester, sodium salt, (2S-cis)- (9CI) (CA INDEX NAME)



● Na

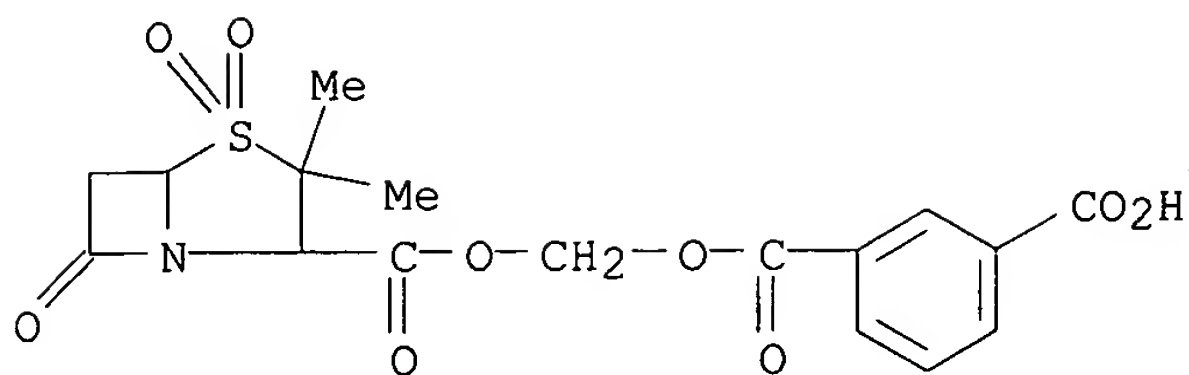
IT 87343-51-9P

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(preparation and reaction of, with ampicillin iodomethyl ester)

RN 87343-51-9 CAPLUS

CN 1,3-Benzenedicarboxylic acid, mono[[[(3,3-dimethyl-4,4-dioxido-7-oxo-4-thia-1-azabicyclo[3.2.0]hept-2-yl)carbonyl]oxy)methyl] ester, sodium salt, (2S-cis)- (9CI) (CA INDEX NAME)



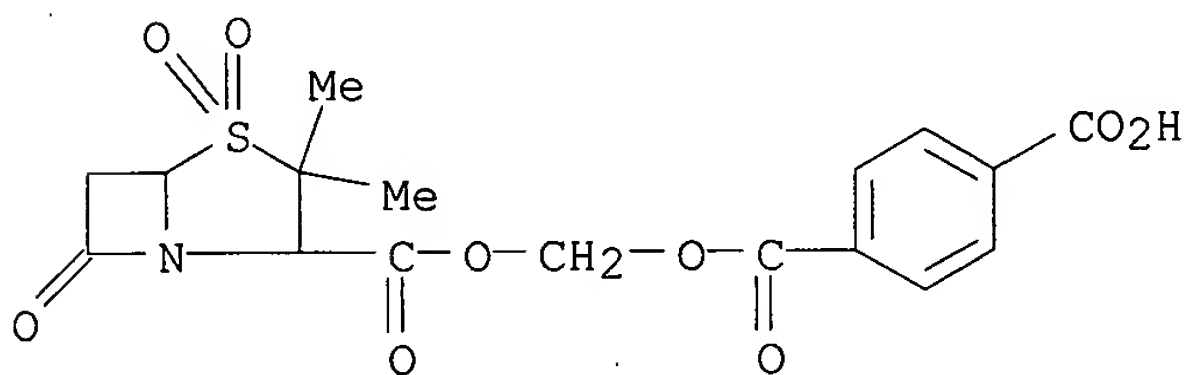
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IT 87343-45-1 87352-92-9 87503-35-3

RL: RCT (Reactant); RACT (Reactant or reagent)  
(reaction of, with acetoacetate)

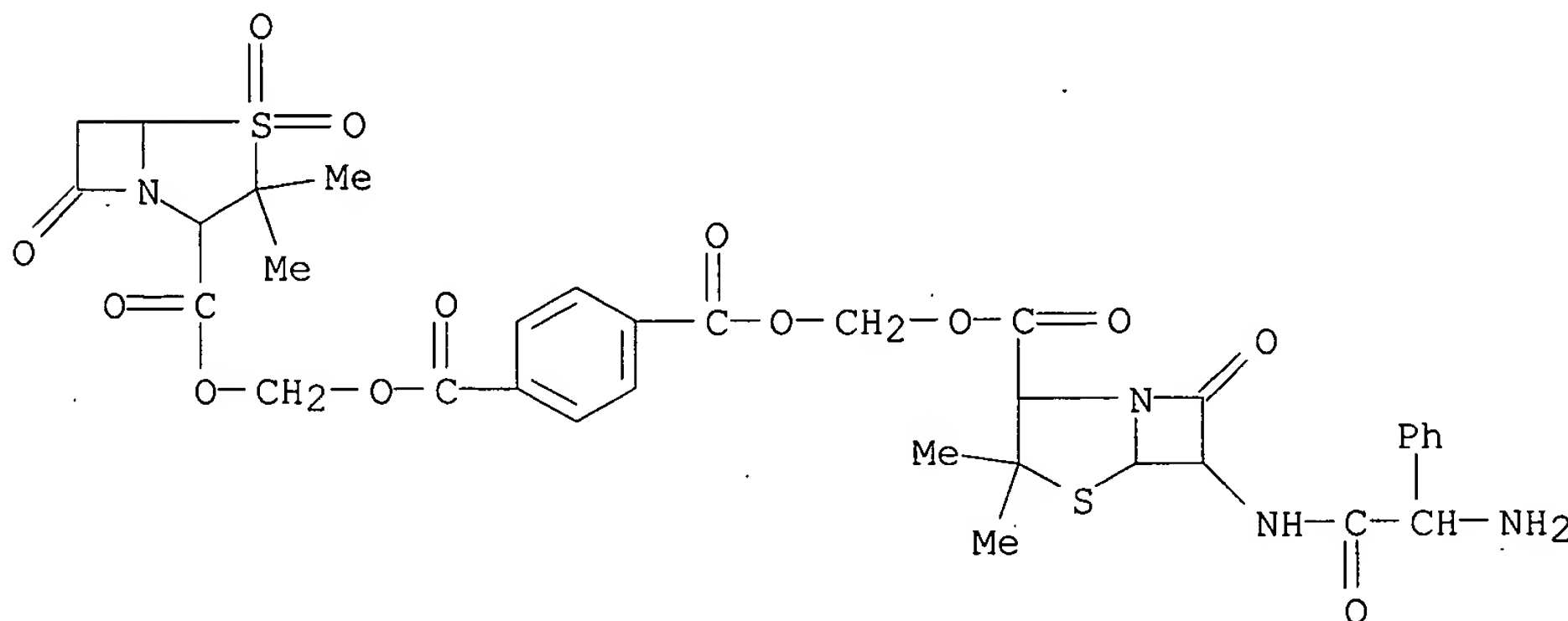
RN 87343-45-1 CAPLUS

CN 1,4-Benzenedicarboxylic acid, mono[[[(3,3-dimethyl-4,4-dioxido-7-oxo-4-thia-1-azabicyclo[3.2.0]hept-2-yl)carbonyl]oxy]methyl] ester, (2S-cis)-(9CI) (CA INDEX NAME)



RN 87352-92-9 CAPLUS

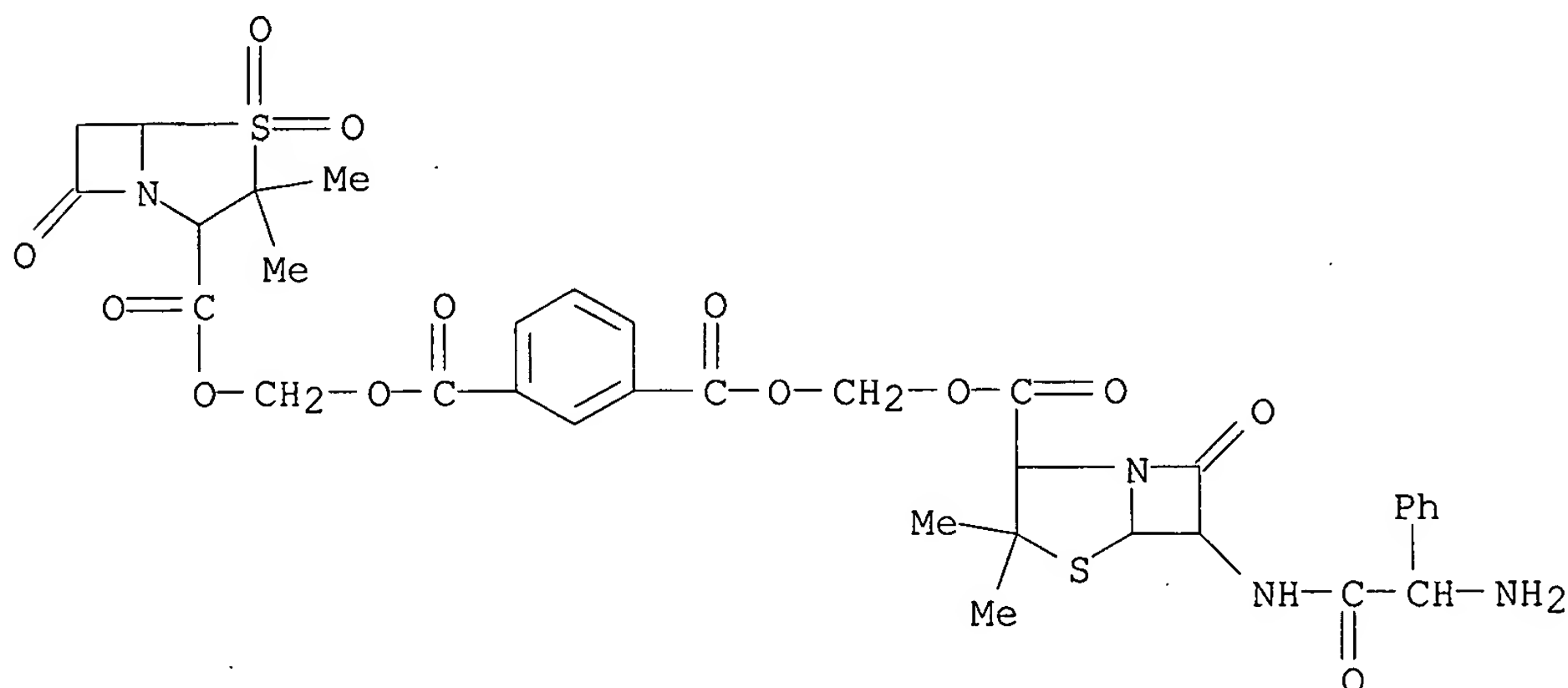
CN 1,4-Benzenedicarboxylic acid, [[[[6-[(aminophenylacetyl)amino]-3,3-dimethyl-7-oxo-4-thia-1-azabicyclo[3.2.0]hept-2-yl]carbonyl]oxy]methyl [[[(3,3-dimethyl-4,4-dioxido-7-oxo-4-thia-1-azabicyclo[3.2.0]hept-2-yl)carbonyl]oxy]methyl ester, monohydrochloride, [2S-[2α(2R\*,5S\*),5α,6β(S\*)]]- (9CI) (CA INDEX NAME)



● HCl

RN 87503-35-3 CAPLUS

CN 1,3-Benzenedicarboxylic acid, [[[6-[(aminophenylacetyl)amino]-3,3-dimethyl-7-oxo-4-thia-1-azabicyclo[3.2.0]hept-2-yl]carbonyl]oxy]methyl  
 [[(3,3-dimethyl-4,4-dioxido-7-oxo-4-thia-1-azabicyclo[3.2.0]hept-2-yl)carbonyl]oxy]methyl ester, monohydrochloride, [2S-[2 $\alpha$ (2R\*,5S\*),5 $\alpha$ ,6 $\beta$ (S\*)]]- (9CI) (CA INDEX NAME)



● HCl

L13 ANSWER 3 OF 4 CAPLUS COPYRIGHT 2005 ACS on STN

ACCESSION NUMBER: 1984:6194 CAPLUS

DOCUMENT NUMBER: 100:6194

TITLE: 1,1-Alkanediol dicarboxylate linked antibacterial agents

INVENTOR(S): Jasys, Vytautas John; Kellogg, Michael Stephen

PATENT ASSIGNEE(S): Pfizer Inc., USA

SOURCE: Eur. Pat. Appl., 124 pp.

CODEN: EPXXDW

DOCUMENT TYPE: Patent

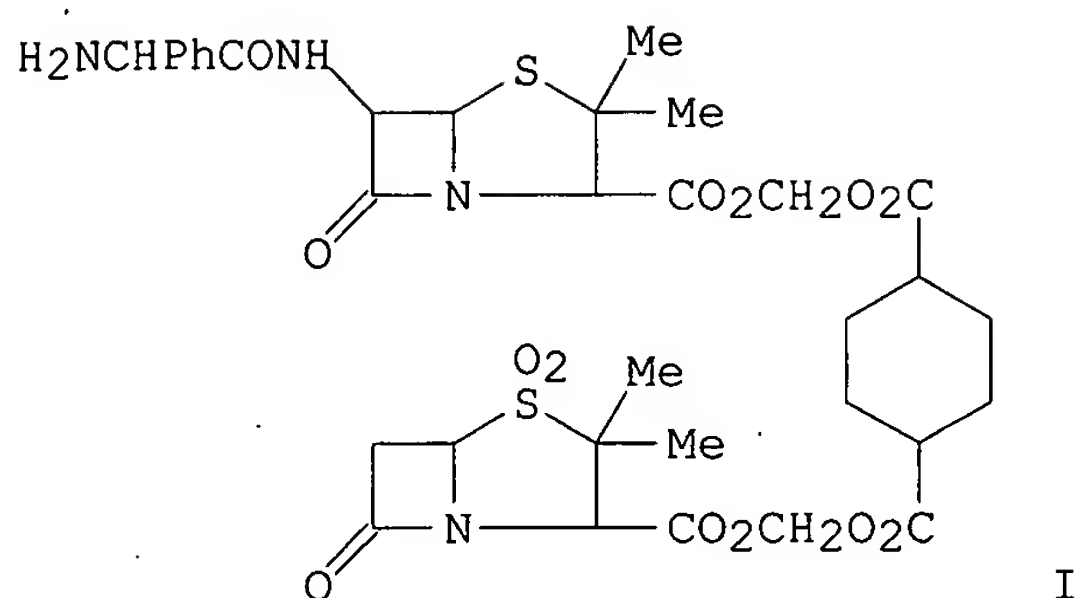
LANGUAGE: English

FAMILY ACC. NUM. COUNT: 2

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
EP 83484	A1	19830713	EP 1982-306683	19821214
EP 83484	B1	19860219		
R: AT, BE, CH, DE, FR, GB, IT, LI, LU, NL, SE				
US 4457924	A	19840703	US 1982-429915	19820930
AT 18051	E	19860315	AT 1982-306683	19821214
PRIORITY APPLN. INFO.:			US 1981-334022	A 19811222
			US 1982-429915	A 19820930
			EP 1982-306683	A 19821214

GI



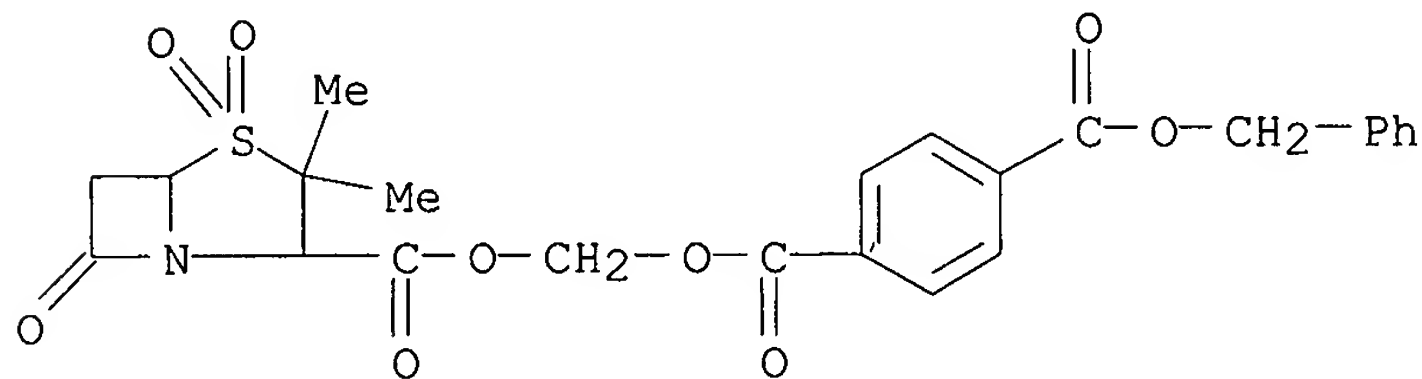
AB Diesters of alkanedicarboxylic acids with penicillin esters and penicillanates, penicillanate dioxides, or hydroxyethyleneoxaazabicycloheptanecarboxylates were prepared. Thus, I was obtained from Na penicillanate dioxide, ampicillin, and K benzyl trans-1,4-cyclohexanedicarboxylate, ClCH<sub>2</sub>I, and BrCH<sub>2</sub>Cl in 10 steps.

IT **87343-43-9P 87343-50-8P**

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)  
(preparation and hydrogenolysis of)

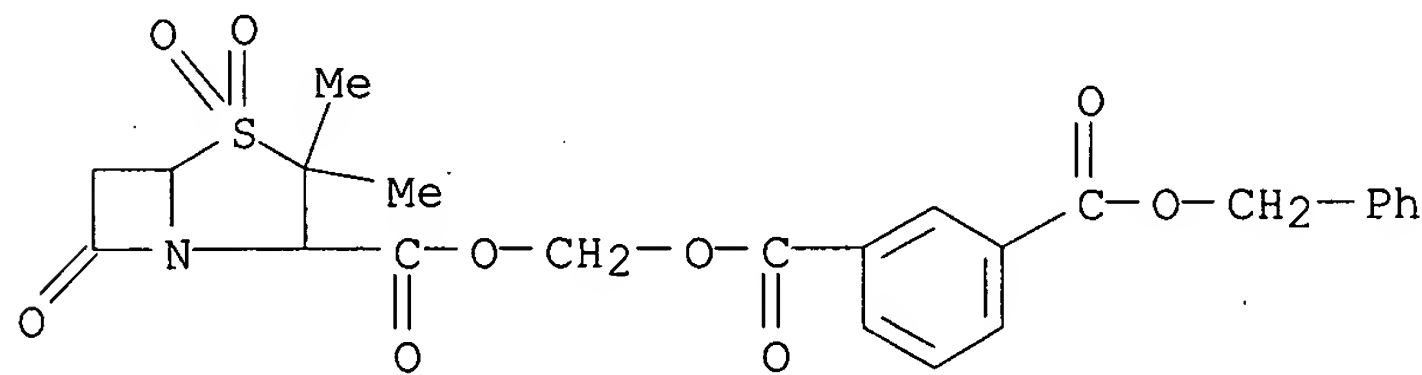
RN 87343-43-9 CAPLUS

CN 1,4-Benzenedicarboxylic acid, [[[3,3-dimethyl-4,4-dioxido-7-oxo-4-thia-1-azabicyclo[3.2.0]hept-2-yl)carbonyl]oxy]methyl phenylmethyl ester, (2S-cis)- (9CI) (CA INDEX NAME)



RN 87343-50-8 CAPLUS

CN 1,3-Benzenedicarboxylic acid, [[[3,3-dimethyl-4,4-dioxido-7-oxo-4-thia-1-azabicyclo[3.2.0]hept-2-yl)carbonyl]oxy]methyl phenylmethyl ester, (2S-cis)- (9CI) (CA INDEX NAME)



IT **87352-91-8P 87352-93-0P**

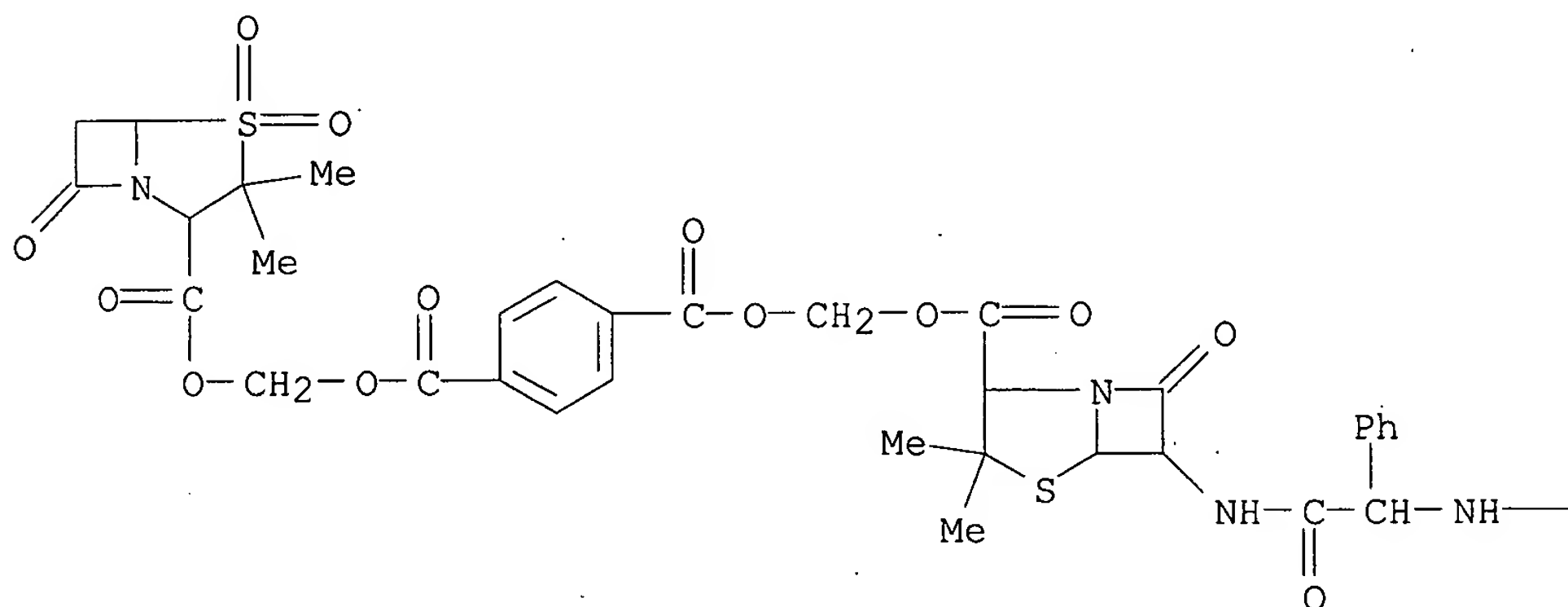
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)  
(preparation and hydrolysis of)

RN 87352-91-8 CAPLUS

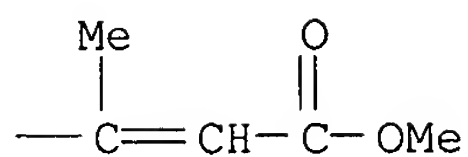


CN 1,4-Benzenedicarboxylic acid, [[(3,3-dimethyl-4,4-dioxido-7-oxo-4-thia-1-azabicyclo[3.2.0]hept-2-yl)carbonyl]oxy]methyl [[[6-[[[(3-methoxy-1-methyl-3-oxo-1-propenyl)amino]phenylacetyl]amino]-3,3-dimethyl-7-oxo-4-thia-1-azabicyclo[3.2.0]hept-2-yl]carbonyl]oxy]methyl ester, [2S-[2 $\alpha$ (2R\*,5S\*),5 $\alpha$ ,6 $\beta$ (S\*)]]- (9CI) (CA INDEX NAME)

PAGE 1-A



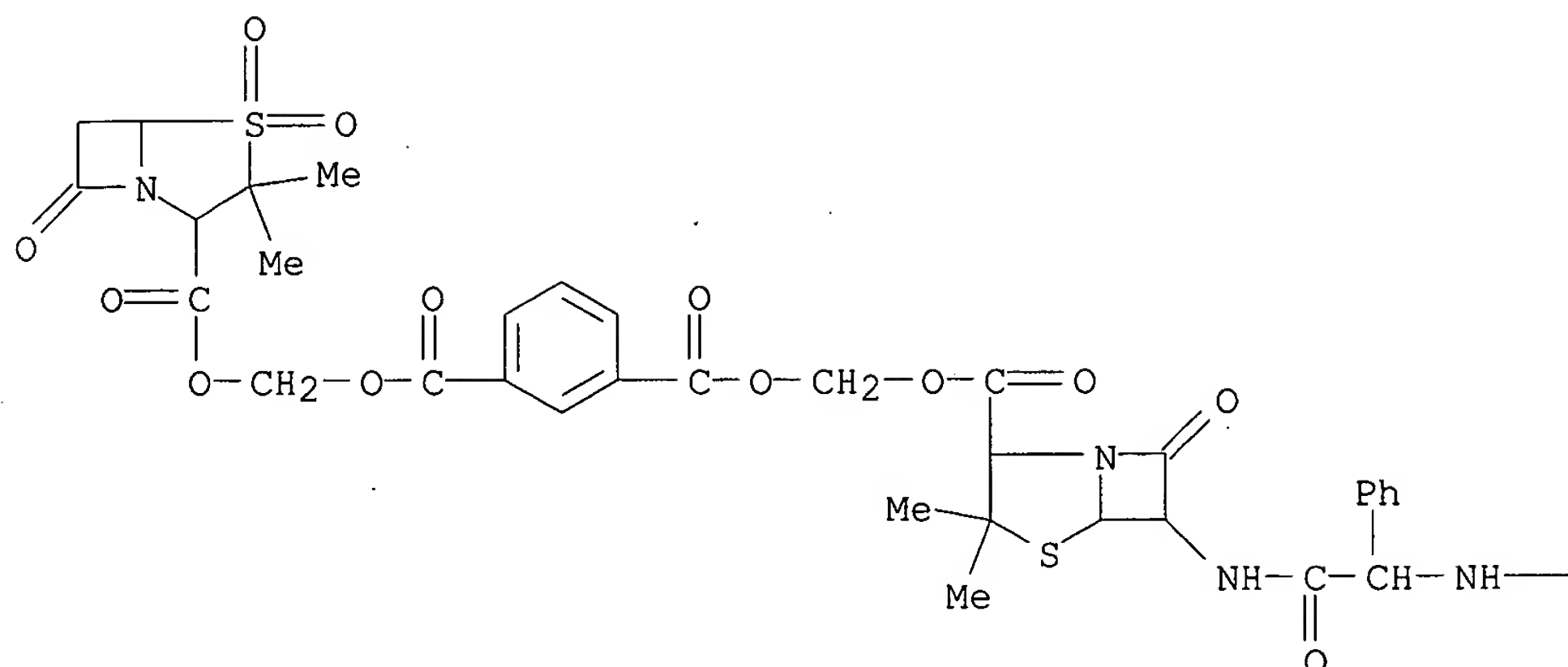
PAGE 1-B



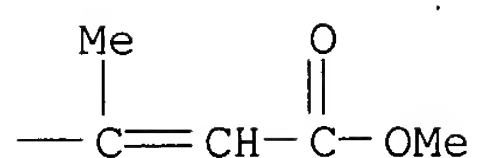
RN 87352-93-0 CAPLUS

CN 1,3-Benzenedicarboxylic acid, [[(3,3-dimethyl-4,4-dioxido-7-oxo-4-thia-1-azabicyclo[3.2.0]hept-2-yl)carbonyl]oxy]methyl [[[6-[[[(3-methoxy-1-methyl-3-oxo-1-propenyl)amino]phenylacetyl]amino]-3,3-dimethyl-7-oxo-4-thia-1-azabicyclo[3.2.0]hept-2-yl]carbonyl]oxy]methyl ester, [2S-[2 $\alpha$ (2R\*,5S\*),5 $\alpha$ ,6 $\beta$ (S\*)]]- (9CI) (CA INDEX NAME)

PAGE 1-A



PAGE 1-B

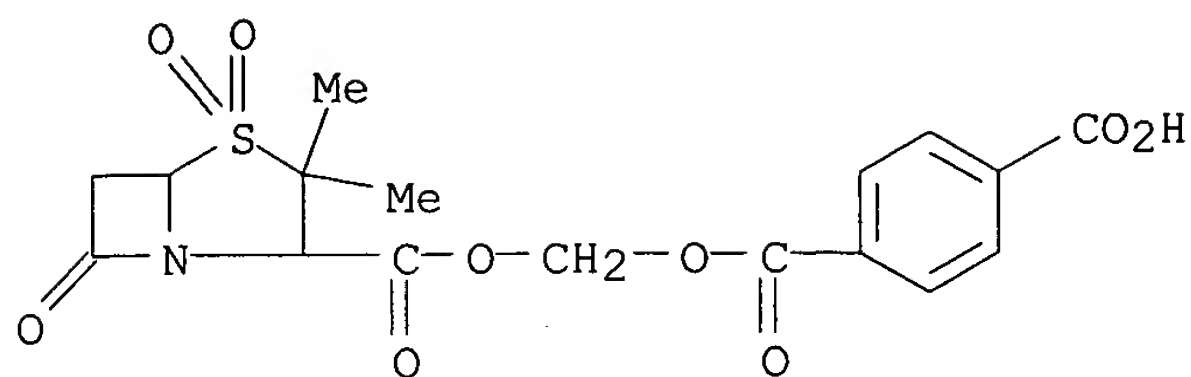


IT 87343-44-0P

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT  
(Reactant or reagent)  
(preparation and neutralization of)

RN 87343-44-0 CAPLUS

CN 1,4-Benzenedicarboxylic acid, mono[[[(3,3-dimethyl-4,4-dioxido-7-oxo-4-thia-1-azabicyclo[3.2.0]hept-2-yl)carbonyl]oxy]methyl] ester, sodium salt,  
(2S-cis)- (9CI) (CA INDEX NAME)



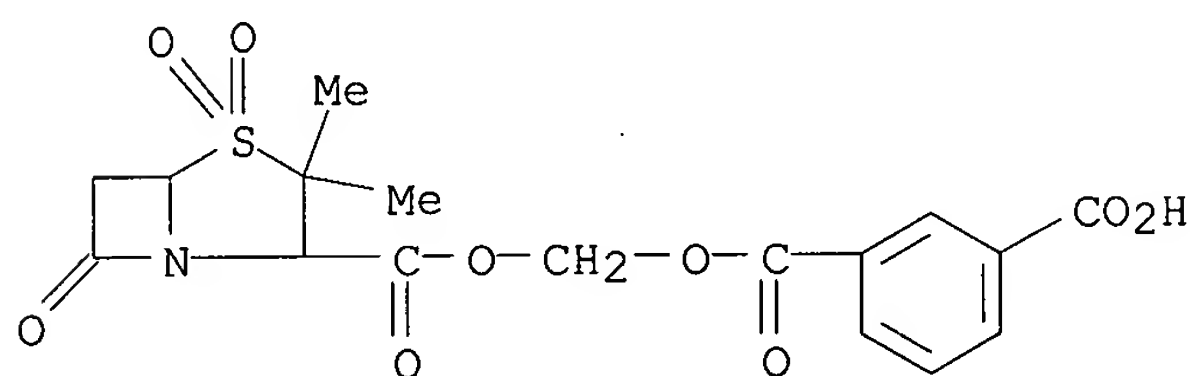
● Na

IT 87343-51-9P

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)  
(preparation and reaction of, with ampicillin iodomethyl ester)

RN 87343-51-9 CAPLUS

CN 1,3-Benzenedicarboxylic acid, mono[[[(3,3-dimethyl-4,4-dioxido-7-oxo-4-thia-1-azabicyclo[3.2.0]hept-2-yl)carbonyl]oxy]methyl] ester, sodium salt, (2S-cis)- (9CI) (CA INDEX NAME)



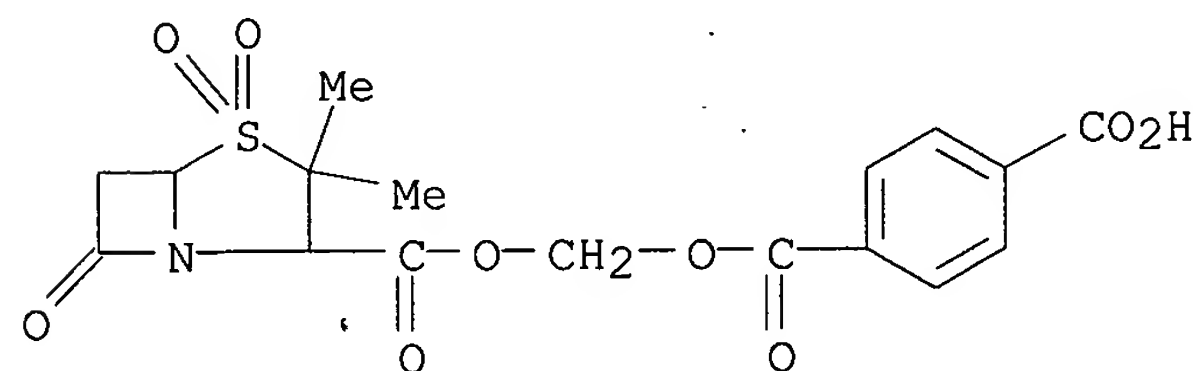
● Na

IT 87343-45-1P 87352-92-9P 87503-35-3P

RL: SPN (Synthetic preparation); PREP (Preparation)  
(preparation of)

RN 87343-45-1 CAPLUS

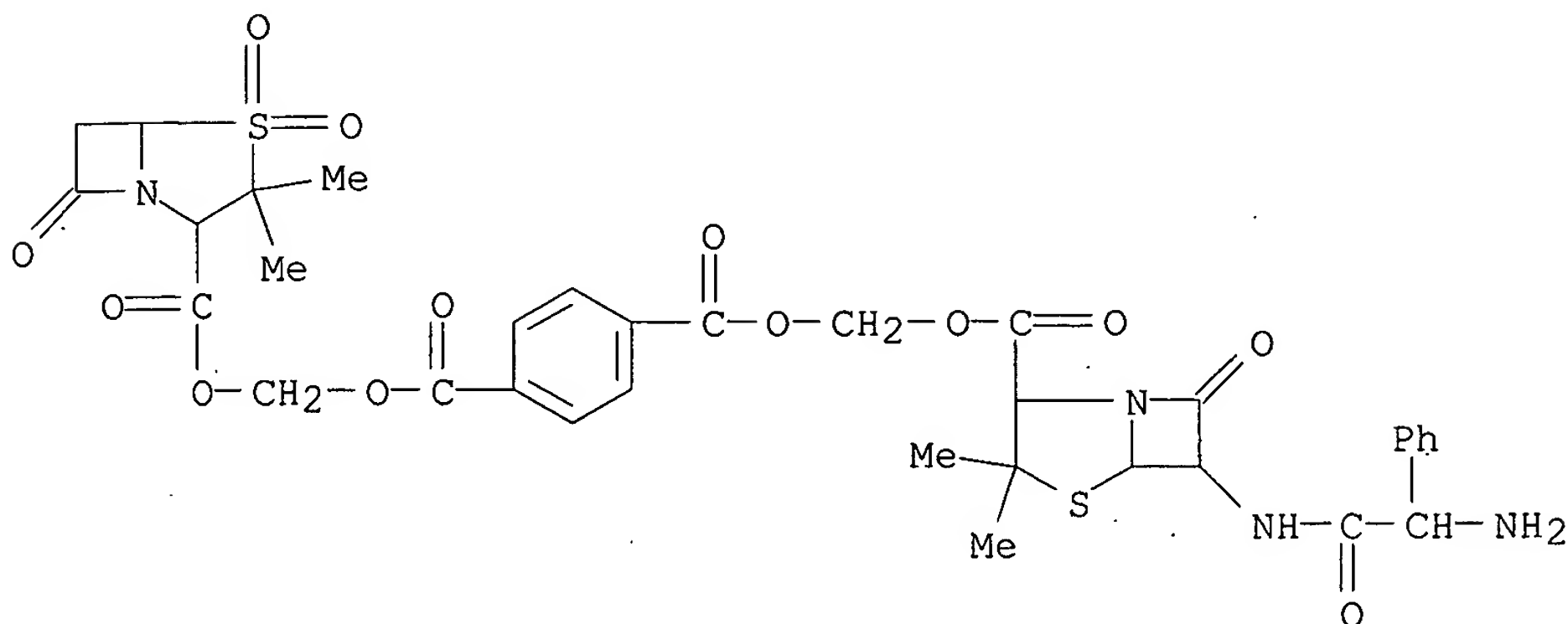
CN 1,4-Benzenedicarboxylic acid, mono[[[(3,3-dimethyl-4,4-dioxido-7-oxo-4-thia-1-azabicyclo[3.2.0]hept-2-yl)carbonyl]oxy]methyl] ester, (2S-cis)- (9CI) (CA INDEX NAME)



RN 87352-92-9 CAPLUS

CN 1,4-Benzenedicarboxylic acid, [[[(6-[(aminophenylacetyl)amino]-3,3-dimethyl-7-oxo-4-thia-1-azabicyclo[3.2.0]hept-2-yl)carbonyl]oxy]methyl] [[(3,3-dimethyl-4,4-dioxido-7-oxo-4-thia-1-azabicyclo[3.2.0]hept-2-

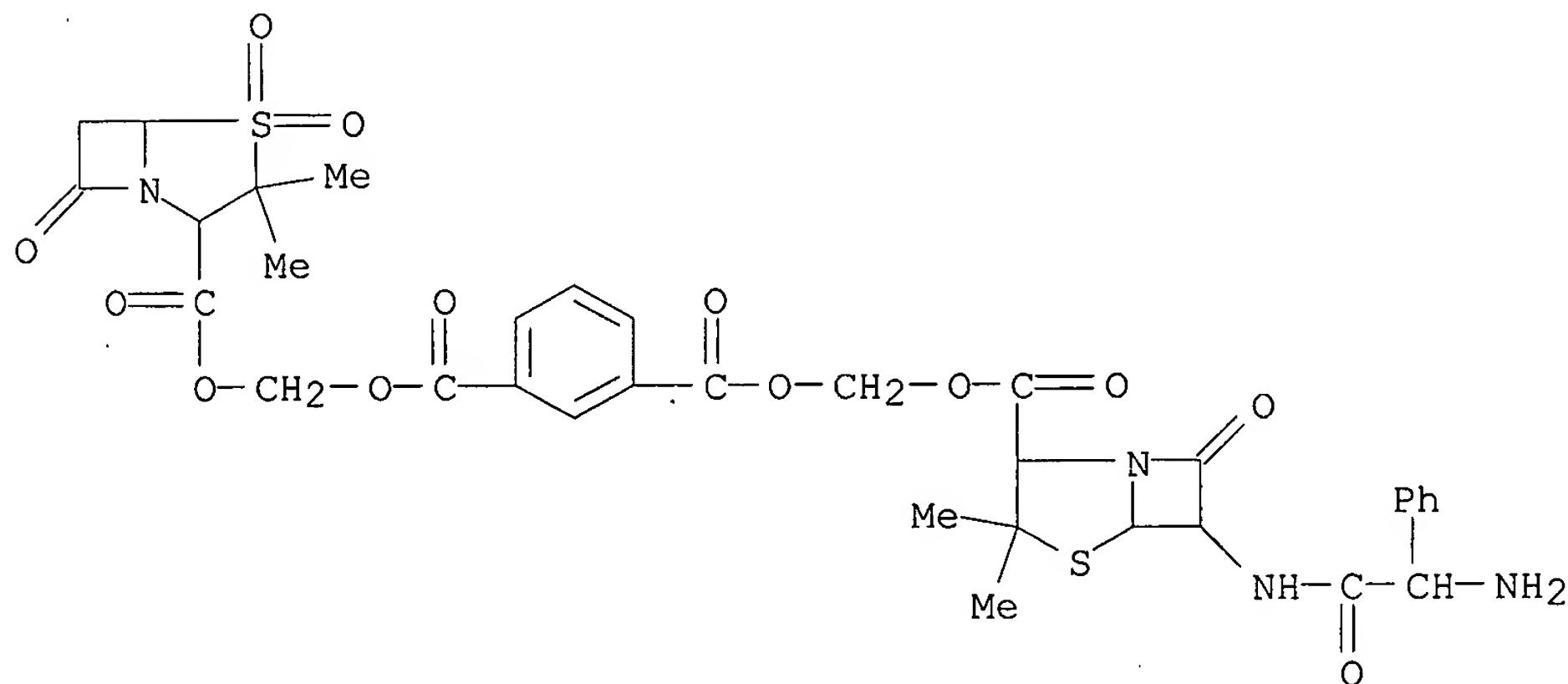
yl)carbonyl]oxy)methyl ester, monohydrochloride, [2S-[2 $\alpha$ (2R\*,5S\*),5 $\alpha$ ,6 $\beta$ (S\*)]]- (9CI) (CA INDEX NAME)



● HCl

RN 87503-35-3 CAPLUS

CN 1,3-Benzenedicarboxylic acid, [[[6-[(aminophenylacetyl)amino]-3,3-dimethyl-7-oxo-4-thia-1-azabicyclo[3.2.0]hept-2-yl]carbonyl]oxy)methyl [[(3,3-dimethyl-4,4-dioxido-7-oxo-4-thia-1-azabicyclo[3.2.0]hept-2-yl)carbonyl]oxy)methyl ester, monohydrochloride, [2S-[2 $\alpha$ (2R\*,5S\*),5 $\alpha$ ,6 $\beta$ (S\*)]]- (9CI) (CA INDEX NAME)



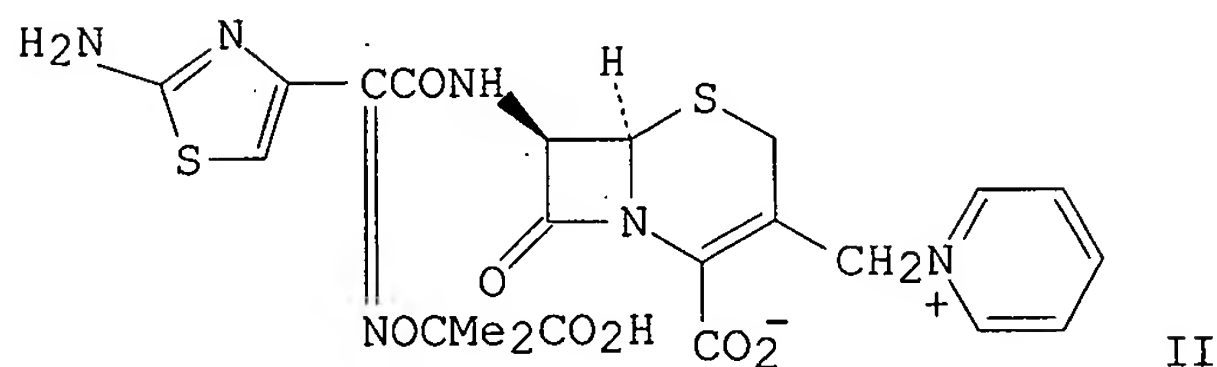
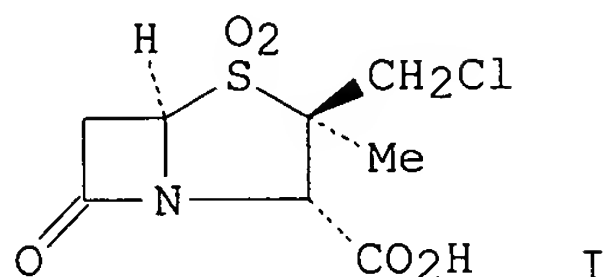
● HCl

L13 ANSWER 4 OF 4 CAPLUS COPYRIGHT 2005 ACS on STN  
 ACCESSION NUMBER: 1983:600528 CAPLUS  
 DOCUMENT NUMBER: 99:200528

TITLE: Treating resistant bacteria including anaerobes  
 INVENTOR(S): Gordon, Maxwell; Pachter, I. Jacob  
 PATENT ASSIGNEE(S): Bristol-Myers Co. , USA  
 SOURCE: U.S., 5 pp.  
 CODEN: USXXAM  
 DOCUMENT TYPE: Patent  
 LANGUAGE: English  
 FAMILY ACC. NUM. COUNT: 1  
 PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
US 4406887	A	19830927	US 1981-310346	19811013
JP 60061528	A2	19850409	JP 1983-165347	19830909
EP 134302	A1	19850320	EP 1983-109144	19830915
EP 134302	B1	19870616		
R: AT, BE, CH, DE, FR, GB, IT, LI, LU, NL, SE				
AT 27772	E	19870715	AT 1983-109144	19830915
PRIORITY APPLN. INFO.:			US 1981-310346	19811013
			EP 1983-109144	A 19830915

GI



AB A synergistic combination of BL-P2013 (I) [79634-05-2] or its salts, as  $\beta$ -lactamase inhibitors, and ceftazidime (II) [72558-82-8] is used to treat bacteroides infection, especially by parenteral administration. Marked synergism was found for I and II against a number of bacteria compared with I or II alone. An injection composition was prepared containing II 500, I K salt [79634-05-2] 500 and Na<sub>2</sub>CO<sub>3</sub> 47 mg.

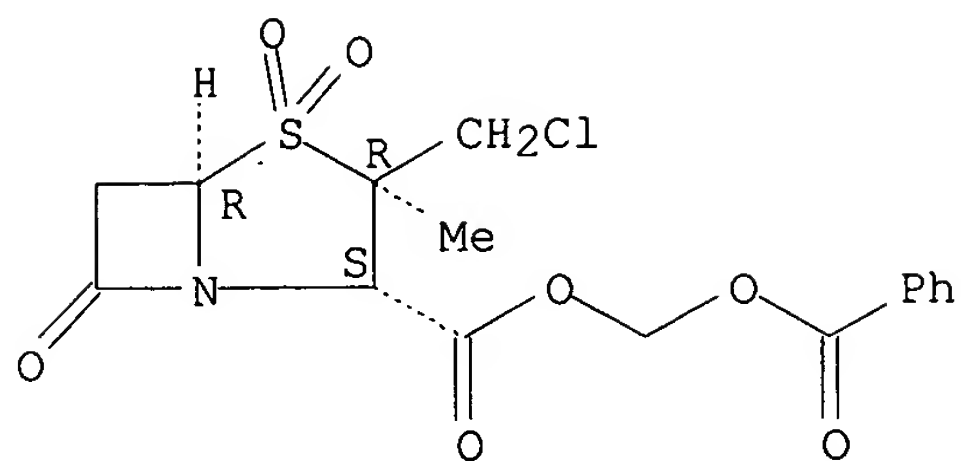
IT 87877-18-7

RL: BIOL (Biological study)  
 (bactericidal compns. containing synergistic combination of ceftazidime with)

RN 87877-18-7 CAPLUS

CN 4-Thia-1-azabicyclo[3.2.0]heptane-2-carboxylic acid, 3-(chloromethyl)-3-methyl-7-oxo-, (benzoyloxy)methyl ester, 4,4-dioxide, [2S-(2 $\alpha$ ,3 $\beta$ ,5 $\alpha$ )]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



=> file beilstein

FILE 'BEILSTEIN' ENTERED AT 15:15:59 ON 18 JUL 2005

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FILE RELOADED ON OCTOBER 20, 2002

FILE LAST UPDATED ON APRIL 21, 2005

FILE COVERS 1771 TO 2004.

\*\*\* FILE CONTAINS 9,218,366 SUBSTANCES \*\*\*

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Reaction data for BEILSTEIN compounds may be displayed  
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(reactions). A substance answer set retrieved after the search  
for a chemical name, a compounds with available reaction  
information by combining with PRE/FA, REA/FA or more generally  
with RX/FA. The BEILSTEIN Registry Number (BRN) is the link  
between a BEILSTEIN compound and belonging reactions. For mo  
detailed reaction searches BRNs can be searched as reaction  
partner BRNs Reactant BRN (RX.RBRN) or Product BRN (RX.PBRN).<<<

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ALL PATENT REFERENCES, OR ALL BABS ACCESSION NUMBERS FOR A  
COMPOUND AT A GLANCE.

=> s L8 full sss

FULL SEARCH INITIATED 15:16:09 FILE 'BEILSTEIN'

FULL SCREEN SEARCH COMPLETED - 570 TO ITERATE

100.0% PROCESSED 570 ITERATIONS  
SEARCH TIME: 00.00.16

0 ANSWERS

L14 0 SEA SSS FUL L8

Berch 10\_648408

07/18/2005

=> d his full

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L1 FILE 'REGISTRY' ENTERED AT 14:25:32 ON 18 JUL 2005  
 L2 STRUCTURE UPLOADED  
 6 SEA SSS SAM L1  
 D SCA

FILE 'STNGUIDE' ENTERED AT 14:29:03 ON 18 JUL 2005

L3 FILE 'REGISTRY' ENTERED AT 14:30:02 ON 18 JUL 2005  
 L4 STRUCTURE UPLOADED  
 5 SEA SSS SAM L3  
 D SCA

L5 FILE 'CAPLUS' ENTERED AT 14:31:38 ON 18 JUL 2005  
 5 SEA ABB=ON PLU=ON L4

L6 FILE 'REGISTRY' ENTERED AT 14:42:20 ON 18 JUL 2005  
 D SCA L4  
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 D L3  
 76 SEA SSS FUL L3

L7 FILE 'CAPLUS' ENTERED AT 14:47:36 ON 18 JUL 2005  
 L8 21 SEA ABB=ON PLU=ON L6  
 STRUCTURE UPLOADED  
 S L8

L9 FILE 'REGISTRY' ENTERED AT 14:56:40 ON 18 JUL 2005  
 1 SEA SSS SAM L8

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 2 SEA ABB=ON PLU=ON L9

L11 FILE 'REGISTRY' ENTERED AT 14:56:59 ON 18 JUL 2005  
 L12 1 SEA SUB=L6 SSS SAM L8  
 17 SEA SUB=L6 SSS FUL L8

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 4 SEA ABB=ON PLU=ON L12

FILE 'REGISTRY' ENTERED AT 14:59:36 ON 18 JUL 2005

FILE 'REGISTRY' ENTERED AT 15:12:34 ON 18 JUL 2005

FILE 'CAPLUS' ENTERED AT 15:12:42 ON 18 JUL 2005  
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L14 FILE 'BEILSTEIN' ENTERED AT 15:15:59 ON 18 JUL 2005  
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FILE 'STNGUIDE' ENTERED AT 15:17:21 ON 18 JUL 2005

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 59 SEA ABB=ON PLU=ON L6 NOT L12



L16 FILE 'CAPLUS' ENTERED AT 15:25:01 ON 18 JUL 2005  
0 SEA ABB=ON PLU=ON L15 AND L13

FILE 'REGISTRY' ENTERED AT 15:26:11 ON 18 JUL 2005

FILE 'CAPLUS' ENTERED AT 15:28:30 ON 18 JUL 2005

FILE HOME

FILE REGISTRY

Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 17 JUL 2005 HIGHEST RN 855596-49-5

DICTIONARY FILE UPDATES: 17 JUL 2005 HIGHEST RN 855596-49-5

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH JANUARY 18, 2005

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\*\*\*\*\*

\*  
\* The CA roles and document type information have been removed from \*  
\* the IDE default display format and the ED field has been added, \*  
\* effective March 20, 2005. A new display format, IDERL, is now \*  
\* available and contains the CA role and document type information. \*  
\*

\*\*\*\*\*

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. For more information enter HELP PROP at an arrow prompt in the file or refer to the file summary sheet on the web at:  
<http://www.cas.org/ONLINE/DBSS/registryss.html>

FILE STNGUIDE

FILE CONTAINS CURRENT INFORMATION.

LAST RELOADED: Jul 15, 2005 (20050715/UP).

FILE CAPLUS

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FILE COVERS 1907 - 18 Jul 2005 VOL 143 ISS 4

FILE LAST UPDATED: 17 Jul 2005 (20050717/ED)

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This file contains CAS Registry Numbers for easy and accurate substance identification.

FILE BEILSTEIN

FILE RELOADED ON OCTOBER 20, 2002

FILE LAST UPDATED ON APRIL 21, 2005

FILE COVERS 1771 TO 2004.

**FILE CONTAINS 9,218,366 SUBSTANCES**

>>>PLEASE NOTE: Reaction Data and substance data are stored in separate documents and can not be searched together in one query. Reaction data for BEILSTEIN compounds may be displayed immediately with the display codes PRE (preparations) and REA (reactions). A substance answer set retrieved after the search for a chemical name, a compounds with available reaction information by combining with PRE/FA, REA/FA or more generally with RX/FA. The BEILSTEIN Registry Number (BRN) is the link between a BEILSTEIN compound and belonging reactions. For more detailed reaction searches BRNs can be searched as reaction partner BRNs Reactant BRN (RX.RBRN) or Product BRN (RX.PBRN).<<<

>>> FOR SEARCHING PREPARATIONS SEE HELP PRE <<<

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